C:\Program Files\Stnexp\Queries\10-676-391-02.str

chain nodes :

exact bonds :

normalized bonds :

12-41 21-25 30-40

35-79

```
27
                        26
       16
          17
               18
                   25
    85
        86
ring nodes :
                                                                       23
                                                                            24
                                                                                28
                                                                                    29
                                                                                        30
                                                                                             31
                                              14
                                                  15
                                                       19
                                                           20
                                                               21
                                                                   22
                5
                          8
                             10
                                 11
                                      12
                                          13
                                                                   50
                                                                        51
                                                                            52
                                                                                53
                                                                                    54
                                                                                             56
                                              45
                                                  46
                                                       47
                                                           48
                                                               49
                    36
                         37
                             38
                                  39
                                      43
                                          44
    32 -33
           34
                35
                                                           73
                                                               74
                                                                   75
                                      68
                                          69
                                              70
                                                  71
                                                       72
                             66
                                 67
    57 58
           59
                60
                    61
                         62
chain bonds :
                                     15-16 16-19 21-25 22-26
                                                                           30-40 35-79
                                                                  23-27
    4-9 9-10
               12-41 13-17 14-18
                                 52-63 54-65 61-81, 67-80 70-78 73-77 75-76
    42-84 43-83 45-82 50-64
    84-86
ring bonds :
                                                          8-28 10-11
                          3-4 3-8 4-5
                                                                        10-15 11-12
                    2-3
                                         5-6
                                               5-7 7-29
    1-2 1-6 1-68
                                                               28-30
                                                                       28-34
                                                                              29-35
                                                                                      29-39
                                  20-21
                                         21-22
                                                22-23
                                                        23-24
                  19-20<sup>.</sup>
                          19-24
           14-15
    13-14
                                                               37-75
                                                                       38-39
                                                                              43-44
                                  33-34
                                         35-36
                                                36-37
                                                        37-38
                  32-33
                          32-43
    30-31
           31-32
                                                                              52-53
                                                                                      54-55
                                                               50-51
                                                                       51-52
           45-46
                                 48-49
                                         48-53
                                                49-50
                                                        49-55
    44-62
                  46-47
                          47-48
                                                                              67-68
                                                        60-66
                                                               61-62
                                                                       66-67
           55-56
                   56-57
                          57-58
                                 57-60
                                         58-59
                                                60-61
    54-59
                  71-72
                                  73-74
                                         74-75
                          72 - 73
    69-70
           70-71
exact/norm bonds :
                                 8-28 9-10 10-11 10-15
                                                                  12-13
                                                                           13-14
                                                                                  13-17
                                                                                          14-15
                                                           11-12
         3-8 4-9
                     5-7
                          7-29
    1-68
                                                                              23-27
                                                                                      32-43
                                                               22-26
                                                                       23-24
          15-16
                          19-20
                                 19-24
                                         20-21
                                                21-22
                                                        22-23
    14-18
                  16-19
                                                44-62
                                                        45-46
                                                               45-82
                                                                       46-47
                                                                              47-48
                                                                                      49-55
           41-42
                  42-84
                          43-44
                                  43-83
                                         44-45
    37-75
                                                               66-67
                                                                       67-68
                                                                              67-80
                                                                                      68-69
                                                61-62
                                                        61-81
                                 60-61
                                         60-66
                   54-65
                          57-60
    50-64
           52-63
                                                                       84-85
                                                                              84-86
                                                               75-76
           70-71
                                  72-73
                                         73-74
                                                73-77
                                                        74-75
    69-70
                   70-78
                          71-72
```

40

41

42

63

64

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78

77

80

79

81

82

83

84

1-2 1-6 2-3 3-4 4-5 5-6 28-30 28-34 29-35 29-39 30-31 31-32 32-33 33-34 35-36 36-37 37-38 38-39 48-49 48-53 49-50 50-51 51-52 52-53 54-55 54-59 55-56 56-57 57-58 58-59

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS 26:CLASS 27:CLASS 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom 56:Atom 57:Atom 58:Atom 59:Atom 60:Atom 61:Atom 62:Atom 63:CLASS 64:CLASS 65:CLASS 66:Atom 67:Atom 68:Atom 69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS 78:CLASS 79:CLASS 80:CLASS 81:CLASS 82:CLASS 83:CLASS 84:CLASS 85:CLASS 86:CLASS

12-41 21-25 30-40

normalized bonds :

35-36

35-79

```
chain nodes :
                                                                                         84
                                                                             81
                                                                                 82
                                                                                     83
                                                        76
                                                            77
                                                                78
                                                                     79
                                                                         80
    9
      16
          17
              18
                   25
                       26
                           27
                               40
                                    41
                                       42 63
                                                64
                                                    65
    85
ring nodes :
                                                                                      30
                                                                                           31
                                                         20
                                                             21
                                                                 22
                                                                     23
                                                                          24
                                                                              28
                                                                                  29
                            10
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                                         13
                                             14
                                                 15
                                                     19
    1 2 3 4
               5
                   6 7
                                                                             53
                                                                                  54
                                                     47
                                                         48
                                                             49
                                                                  50
                                                                      51
                                                                          52
    32 33 34 35 36
                        37
                            38
                                39
                                     43
                                         44
                                             45
                                                 46
                                     68
                                         69
                                             70
                                                 71
                                                     72
                                                         73
                                                             74
                                                                  75
    57 58 59 60 61
                        62
                            66
                                67
chain bonds :
    4-9 9-10 12-41 13-17 14-18 15-16 16-19 21-25 22-26
                                                                 23-27 30-40 35-79 41-42
                                50-64 52-63 54-65 61-81 67-80 70-78 73-77 75-76
    42-84 42-85 43-83 45-82
                         3-4 3-8 4-5 5-6 5-7 7-29
                                                         8-28 10-11 10-15
                                                                             11-12 12-13
    1-2 1-6 1-68 2-3
                                                                                   29-39
                                                             28-30
                                                                     28-34
                                                                            29-35
                                              22-23
                                                      23-24
    13-14 14-15
                  19-20
                         19-24
                                20-21
                                        21-22
                                               36-37
                                                      37-38
                                                             37-75
                                                                     38-39
                                                                            43-44
                         32-43
                                33-34
                                        35-36
    30-31 -31-32
                  32-33
                                        48-53
                                               49-50
                                                      49-55
                                                             50-51
                                                                     51-52
                                                                            52-53
                                                                                   54-55
                         47-48
                                48-49
    44-62
          45-46
                 46-47
                                                                            67-68
                                                                                   68-69
                                                             61-62
                                                                     66-67
                                57-60
                                        58-59
                                               60-61
                                                      60-66
    54-59
           55-56
                 56-57
                         57-58
                                73-74
                                        74-75
    69-70 70-71
                 71-72
                         72-73
exact/norm bonds :
                                8-28 \quad 9-10 \quad 10-11 \quad 10-15 \quad 11-12 \quad 12-13 \quad 13-14
                                                                                13-17 14-15
                         7-29
    1-68 3-8 4-9 5-7
                                        20-21
                                               21-22
                                                      22-23
                                                              22-26
                                                                     23-24
                                                                            23-27
                                                                                   32-43
                         19-20
                                19-24
    14-18 15-16
                 16-19
                                                                                   47-48
                                                              45-46
                                                                     45-82
                                                                            46-47
                                               44-45
                                                      44-62
    37-75
          41-42
                 42-84
                         42-85
                                43-44
                                        43-83
                  52-63
                         54-65
                                 57-60
                                       60-61
                                               60-66
                                                      61-62
                                                              61-81
                                                                     66-67
                                                                            67-68
    49-55
           50-64
                                71-72
                                       72-73
                                               73-74
                                                      73-77
                                                              74-75
                                                                     75-76
           69-70
                  70-71
                         70-78
    68-69
exact bonds :
```

1-2 1-6 2-3 3-4 4-5 5-6 28-30 28-34 29-35 29-39 30-31 31-32 32-33 33-34

36-37 37-38 38-39 48-49 48-53 49-50 50-51 51-52 52-53 54-55 54-59 55-56 56-57 57-58 58-59

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS 26:CLASS 27:CLASS 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom 56:Atom 57:Atom 58:Atom 59:Atom 60:Atom 61:Atom 62:Atom 63:CLASS 64:CLASS 65:CLASS 66:Atom 67:Atom 68:Atom 69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS 78:CLASS 79:CLASS 80:CLASS 81:CLASS 82:CLASS 83:CLASS 84:CLASS 85:CLASS

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NEWS
        DEC 18
                with preparation role
                CA/CAplus patent kind codes updated
        DEC 18
NEWS
                MARPAT to CA/CAplus accession number crossover limit increased
        DEC 18
NEWS 5
                 to 50,000
                MEDLINE updated in preparation for 2007 reload
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        DEC 18
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        DEC 27
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NEWS 13 JAN 22
                CA/CAplus enhanced with patent applications from India
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                PHAR reloaded with new search and display fields
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                 multiple databases
                PATDPASPC enhanced with Drug Approval numbers
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                KOREAPAT enhanced with IPC 8 features and functionality
NEWS 18 FEB 23
NEWS 19 FEB 26
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                EMBASE enhanced with Clinical Trial Number field
NEWS 20 FEB 26
                TOXCENTER enhanced with reloaded MEDLINE
NEWS 21 FEB 26
                IFICDB/IFIPAT/IFIUDB reloaded with enhancements
NEWS 22 FEB 26
                CAS Registry Number crossover limit increased from 10,000
NEWS 23 FEB 26
                 to 300,000 in multiple databases
                WPIDS/WPIX enhanced with new FRAGHITSTR display format
NEWS 24 MAR 15
NEWS 25 MAR 16 CASREACT coverage extended
NEWS 26 MAR 20 MARPAT now updated daily
NEWS 27 MAR 22 LWPI reloaded
NEWS 28 MAR 30
                RDISCLOSURE reloaded with enhancements
        MAR 30
                 INPADOCDB will replace INPADOC on STN
NEWS 29
                JICST-EPLUS removed from database clusters and STN
NEWS 30 APR 02
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              MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
              AND CURRENT DISCOVER FILE IS DATED 25 SEPTEMBER 2006.
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NEWS LOGIN
              For general information regarding STN implementation of IPC 8
NEWS IPC8
              X.25 communication option no longer available
NEWS X25
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=>
Uploading C:\Program Files\Stnexp\Queries\10-676-391-01.str

```
chain nodes :
 9 16 17 18 25 26 27 40 41 42 63 64 65 76 77 78 79 80
                                                                                81
                                                                                    82
 84 85
 ring nodes :
 1 2 3 4 5 6 7 8 10 11
                                     12
                                         13
                                              14 15
                                                       19
                                                           20
                                                                21
                                                                     22
                                                                         23
                                                                              24
                                                                                  28
                                                                                       29
                                                                                           30
 31 32 33 34 35 36 37
                               38
                                     39
                                         43
                                              44
                                                  45
                                                       46
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                                                                     49
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                                                                              51
                                                                                  52 53 54
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                                                                72
                                                                     73
                                                                         74
                                                                              75
 55 56 57 58 59 60 61
                                62
                                     66
                                         67
                                              68
 chain bonds :
 4-9 9-10 12-41 13-17 14-18 15-16 16-19 21-25 22-26 23-27 30-40 35-79
 41-42 42-84 42-85 43-83 45-82 50-64 52-63 54-65 61-81 67-80 70-78 73-77
 75-76
 ring bonds :
                        3-4 3-8 4-5 5-6 5-7 7-29 8-28 10-11 10-15 11-12
 1-2 1-6 1-68 2-3
                        19-20 19-24 20-21 21-22 22-23 23-24 28-30 28-34 29-35
 12-13 13-14 14-15
                                                        36-37 37-38
                                                                       37-75 38-39 43-44
                                                35-36
 29-39 30-31 31-32
                        32-33 32-43
                                        33-34
                                                        49-50 49-55 50-51 51-52 52-53
 44-45 44-62 45-46
                        46-47 47-48
                                        48-49
                                                48-53
                                                        60-61 60-66 61-62 66-67 67-68
 54-55 54-59 55-56
                        56-57 57-58
                                        57-60 58-59
 68-69 69-70 70-71
                        71-72 72-73 73-74
                                                74-75
 exact/norm bonds :
                         7-29 8-28 9-10 10-11 10-15 11-12 12-13 13-14 13-17
. 1-68 3-8 4-9 5-7
                        16-19 19-20 19-24 20-21 21-22 22-23 22-26
42-84 42-85 43-44 43-83 44-45 44-62 45-46
 14-15 14-18 15-16
                                                                                23-24
                                                                                        23-27

    42-84
    42-85
    43-44
    43-83
    44-45
    44-62
    45-46
    45-82

    52-63
    54-65
    57-60
    60-61
    60-66
    61-62
    61-81
    66-67

    70-71
    70-78
    71-72
    72-73
    73-74
    73-77
    74-75
    75-76

 32-43 37-75 41-42
                                                                                45-82 46-47
                                                                                66-67
                                                                                        67-68
 47-48 49-55 50-64
 67-80 68-69 69-70
 exact bonds :
 12-41 21-25 30-40 35-79
 normalized bonds :
 1-2 \quad 1-6 \quad 2-3 \quad 3-4 \quad 4-5 \quad 5-6 \quad 28-30 \quad 28-34 \quad 29-35 \quad 29-39 \quad 30-31 \quad 31-32 \quad 32-33
 33-34 35-36 36-37 37-38 38-39 48-49 48-53 49-50 50-51 51-52 52-53 54-55 54-59 55-56 56-57 57-58 58-59
```

```
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS 26:CLASS 27:CLASS 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom 56:Atom 57:Atom 58:Atom 59:Atom 60:Atom 61:Atom 62:Atom 63:CLASS 64:CLASS 65:CLASS 66:Atom 75:Atom 68:Atom 69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS 78:CLASS 79:CLASS 80:CLASS 81:CLASS 82:CLASS 83:CLASS 84:CLASS 85:CLASS
```

L1 STRUCTURE UPLOADED

=> d l1 L1 HAS NO ANSWERS

L1

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam

SAMPLE SEARCH INITIATED 15:56:28 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 0 TO 0 PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 sss ful FULL SEARCH INITIATED 15:56:35 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 20 TO ITERATE

. 100.0% PROCESSED 20 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

L3 0 SEA SSS FUL L1

Uploading C:\Program Files\Stnexp\Queries\vancomycin core #3.str

```
chain nodes :
                                          65
                                             76
                                                77
                                                     78
                                                        79
                                                            80
                                                                81
                                                                   82
                                                                       83
9 16 17 18 25 26 27 40 41
                              42
                                  63
                                      64
ring nodes :
                                              20
                                                  21
                                                     22
                                                         23
                                                             24
                                                                28
                                                                    29
                                                                        30
1 2 3 4 5 6 7 8
                     10
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                                   14 15 19
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                            12
31 32 33 34 35 36
                    37
                         38
                            39
                               43
                                   44
                                      45 46
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55 56 57 58 59 60 61
                        62
                            66
                               67
                                   68
                                      69
                                          70
                                              71
                                                  72
                                                      73
                                                         74
                                                             75
chain bonds :
4-9 9-10 12-41 13-17 14-18 15-16 16-19 21-25 22-26 23-27 30-40 35-79
41-42 43-83 45-82 50-64 52-63 54-65 61-81 67-80 70-78 73-77 75-76
ring bonds :
                  3-4 3-8 4-5 5-6 5-7 7-29 8-28 10-11 10-15 11-12
1-2 1-6 1-68 2-3
                              20-21
                                    21-22 22-23 23-24 28-30 28-34 29-35
12-13 13-14 14-15
                  19-20 19-24
                                                        37-75
                                                               38-39 43-44
                  32-33
                                            36-37
                                                  37-38
                         32-43
                               33-34
                                     35-36
29-39 30-31 31-32
                                                  49-55 50-51
                                                               51-52 52-53
                                     48-53
                                           49-50
                  46-47
                        47-48
                               48-49
44-45 44-62 45-46
                               57-60
                                     58-59
                                           60-61 60-66 61-62 66-67 67-68
54-55 54-59 55-56
                  56-57
                        57-58
68-69 69-70 70-71 71-72 72-73
                               73-74
                                     74-75
exact/norm bonds :
1-68 3-8 4-9 5-7
                  7-29 8-28 9-10 10-11 10-15 11-12 12-13 13-14 13-17
14-15 14-18 15-16
                  16-19 19-20 19-24 20-21 21-22 22-23
                                                        22-26
                                                               23-24
                                                        46-47 47-48 49-55
                        43-83
32-43 37-75 41-42
                                           45-46
                                                  45-82
                  43-44
                              44-45
                                     44-62
                                                  66-67 67-68 67-80 68-69
50-64 52-63 54-65
                              60-66 61-62
                                           61-81
                  57-60
                        60-61
                                    73-77 74-75
                                                  75-76
69-70 70-71 70-78
                  71-72 72-73 73-74
exact bonds :
12-41 21-25 30-40 35-79
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 28-30 28-34 29-35 29-39 30-31 31-32 32-33
                  37-38 38-39 48-49 48-53 49-50 50-51 51-52 52-53 54-55
33-34 35-36 36-37
54-59 55-56 56-57 57-58 58-59
```

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom

11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS 26:CLASS 27:CLASS 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 33:Atom 33:Atom 37:Atom

38:Atom 39:Atom 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom

Match level :

STRUCTURE UPLOADED

=> d 14

L4

L4 HAS NO ANSWERS

L4 · STI

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=> s 14 sss sam

SAMPLE SEARCH INITIATED 15:57:32 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 53 TO ITERATE

100.0% PROCESSED

53 ITERATIONS

12 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS:

624 TO 1496

PROJECTED ANSWERS:

33 TO 447

L5 12 SEA SSS SAM L4

=> d his

(FILE 'HOME' ENTERED AT 15:55:45 ON 12 APR 2007)

FILE 'REGISTRY' ENTERED AT 15:55:53 ON 12 APR 2007

L1 STRUCTURE UPLOADED

L2 0 S L1 SSS SAM

L3 0 S L1 SSS FUL

L4 STRUCTURE UPLOADED

L5 12 S L4 SSS SAM

=> log hold

COST IN U.S. DOLLARS SINCE FILE TOTAL

FULL ESTIMATED COST ENTRY SESSION 173.00 173.21

SESSION WILL BE HELD FOR 120 MINUTES STN INTERNATIONAL SESSION SUSPENDED AT 15:57:55 ON 12 APR 2007

SIN INTERNATIONAL SESSION SUSPENDED AT 13.37.33 ON 12 Min 200

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssptafxl1600

PASSWORD:

* * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'REGISTRY' AT 16:00:38 ON 12 APR 2007 FILE 'REGISTRY' ENTERED AT 16:00:38 ON 12 APR 2007 COPYRIGHT (C) 2007 American Chemical Society (ACS)

SINCE FILE TOTAL ENTRY SESSION

COST IN U.S. DOLLARS

FULL ESTIMATED COST 173.00 173.21

=> d his

(FILE 'HOME' ENTERED AT 15:55:45 ON 12 APR 2007)

FILE 'REGISTRY' ENTERED AT 15:55:53 ON 12 APR 2007

L1 STRUCTURE UPLOADED

L2 0 S L1 SSS SAM L3 0 S L1 SSS FUL

STRUCTURE UPLOADED

L5 12 S L4 SSS SAM

=>

L4

Uploading C:\Program Files\Stnexp\Queries\10-676-391-02.str

chain nodes : 9 16 17 18 25 26 27 40 41 42 63 64 65 76 77 78 79 80 81 82 83 84 85 86 ring nodes : 1 2 3 4 5 6 7 8 24 28 20 21 22 23 29 10 11 12 13 14 15 19 52 53 54 31 32 33 34 35 36 55 56 57 58 59 60 47 48 49 50 51 37 38 39 43 44 45 46 75 73 61 62 66 67 68 69 70 71 72 chain bonds : 4-9 9-10 12-41 13-17 14-18 15-16 16-19 21-25 22-26 23-27 30-40 35-79 41-42 42-84 43-83 45-82 50-64 52-63 54-65 61-81 67-80 70-78 73-77 75-76 84-85 84-86 ring bonds : 3-4 3-8 4-5 5-6 5-7 7-29 8-28 10-11 10-15 11-12 1-2 1-6 1-68 2-3 20-21 21-22 22-23 23-24 28-30 28-34 29-35 12-13 13-14 14-15 19-20 19-24 36-37 37-38 37-75 38-39 43-44 33-34 35-36 32-43 29-39 30-31 31-32 32-33 49-50 49-55 50-51 51-52 52-53 48-49 48-53 46-47 47-48 44-45 44-62 45-46 60-61 60-66 61-62 66-67 67-68 56-57 57-58 57-60 58-59 54-55 54-59 55-56 68-69 69-70 70-71 71-72 72-73 73-74 74-75 exact/norm bonds :

```
1-68 3-8 4-9 5-7 7-29 8-28 9-10 10-11 10-15 11-12 12-13 13-14 13-17
14-15 14-18 15-16 16-19 19-20 19-24 20-21 21-22 22-23 22-26 23-24 23-27
32-43 37-75 41-42 42-84 43-44 43-83 44-45 44-62 45-46 45-82 46-47 47-48
49-55 50-64 52-63 54-65 57-60 60-61 60-66 61-62 61-81 66-67 67-68 67-80
68-69 69-70 70-71 70-78 71-72 72-73 73-74 73-77 74-75 75-76 84-85 84-86
exact bonds:
12-41 21-25 30-40 35-79
normalized bonds:
1-2 1-6 2-3 3-4 4-5 5-6 28-30 28-34 29-35 29-39 30-31 31-32 32-33
33-34 35-36 36-37 37-38 38-39 48-49 48-53 49-50 50-51 51-52 52-53 54-55
54-59 55-56 56-57 57-58 58-59
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Match level : 1:Atom 2:Atom

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:CLASS 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:CLASS 17:CLASS 18:CLASS 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS 26:CLASS 27:CLASS 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:CLASS 41:CLASS 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom 56:Atom 57:Atom 58:Atom 59:Atom 60:Atom 61:Atom 62:Atom 63:CLASS 64:CLASS 65:CLASS 66:Atom 67:Atom 68:Atom 69:Atom 70:Atom 71:Atom 72:Atom 73:Atom 74:Atom 75:Atom 76:CLASS 77:CLASS 78:CLASS 79:CLASS 80:CLASS 81:CLASS 82:CLASS 83:CLASS 84:CLASS 85:CLASS 86:CLASS

L6 STRUCTURE UPLOADED

=> d 16 L6 HAS NO ANSWERS L6 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> s 16 sss sam
SAMPLE SEARCH INITIATED 16:02:04 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 0 TO ITERATE

100.0% PROCESSED 0 ITERATIONS 0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 0 TO 0 PROJECTED ANSWERS: 0 TO 0

L7 0 SEA SSS SAM L6

=> s 16 sss ful FULL SEARCH INITIATED 16:02:13 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 20 TO ITERATE

100.0% PROCESSED 20 ITERATIONS 13 ANSWERS SEARCH TIME: 00.00.01

L8 13 SEA SSS FUL L6

=> d 18 1-13

L8 ANSWER 1 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN

- RN 370564-22-0 REGISTRY
- ED Entered STN: 17 Nov 2001
- CN Vancomycin, N3''-[[4-[(3,4-dichlorophenyl)methoxy]phenyl]methyl]-, 6'-(2,4,6-trimethylbenzenesulfonate) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C89 H95 Cl4 N9 O27 S
- SR CA
- LC STN Files: CA, CAPLUS, USPAT2

Absolute stereochemistry.

PAGE 1-B

ОН

PAGE 3-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)

ОН

- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 2 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256351-09-4 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, N3'',56-bis[(phenylmethoxy)carbonyl]-, phenylmethyl ester, 6'-(2,4,6-trimethylbenzenesulfonate) (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C98 H103 Cl2 N9 O30 S

SR CA

LC STN Files: CA, CAPLUS, CASREACT, USPATFULL

Absolute stereochemistry.

PAGE 1-A

PAGE 1-B

∕_Bu-i

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 3 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256351-08-3 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, 6'-(2,4,6-trimethylbenzenesulfonate), mono(trifluoroacetate) (salt) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C75 H85 Cl2 N9 O26 S . C2 H F3 O2
- SR CA
- LC STN Files: CA, CAPLUS, CASREACT, USPATFULL

CM 1

CRN 256350-16-0

CMF C75 H85 Cl2 N9 O26 S

Absolute stereochemistry.

ОН

NH₂

CM 2

CRN 76-05-1 CMF C2 H F3 O2

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- ANSWER 4 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN L8
- RN
- 256351-00-5 REGISTRY Entered STN: 17 Feb 2000 ED
- Vancomycin, 26-decarboxy-26-[[(2-methoxy-2-oxoethyl)amino]carbonyl]-, CN 6'-(2,4,6-trimethylbenzenesulfonate) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C78 H90 Cl2 N10 O27 S
- SR CA

LC STN Files: CA, CAPLUS, CASREACT, USPATFULL

Absolute stereochemistry.

PAGE 1-A

PAGE 2-A

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 5 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256350-95-5 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, 26-decarboxy-26-[[(2-methoxy-2-oxoethyl)amino]carbonyl]N3'',56-bis[(2-propenyloxy)carbonyl]-, 6'-(2,4,6trimethylbenzenesulfonate) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH

C86 H98 Cl2 N10 O31 S MF

SR LC STN Files: CA, CAPLUS, CASREACT, USPATFULL

Absolute stereochemistry.

PAGE 1-A

PAGE 1-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 6 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256350-24-0 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, 6'-deoxy-6'-[(2-pyrenylsulfonyl)oxy]-, mono(trifluoroacetate) (salt) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C82 H83 Cl2 N9 O26 S . C2 H F3 O2
- SR CA
- LC STN Files: CA, CAPLUS, CASREACT, USPATFULL

CM 1

CRN 256350-23-9

CMF C82 H83 Cl2 N9 O26 S

Absolute stereochemistry.

PAGE 1-B

PAGE 2-B

PAGE 3-B

CM 2

CRN · 76-05-1 CMF C2 H F3 O2

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L8 ANSWER 7 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN

RN 256350-23-9 REGISTRY

ED Entered STN: 17 Feb 2000

CN Vancomycin, 6'-deoxy-6'-[(2-pyrenylsulfonyl)oxy]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C82 H83 Cl2 N9 O26 S

CI COM

SR CA

Absolute stereochemistry.

PAGE 1-A

PAGE 3-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L8 ANSWER 8 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN

ОН

RN 256350-20-6 REGISTRY

ED Entered STN: 17 Feb 2000

FS STEREOSEARCH

MF C85 H105 Cl2 N9 O26 S . C2 H F3 O2

SR CA

LC STN Files: CA, CAPLUS, USPATFULL

CM 1

CRN 256350-19-3

CMF C85 H105 Cl2 N9 O26 S

Absolute stereochemistry.

ОН

CM 2

CRN 76-05-1 CMF C2 H F3 O2

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 9 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256350-19-3 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, N3''-decyl-6'-deoxy-6'-[[(2,4,6-trimethylphenyl)sulfonyl]oxy]-(9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C85 H105 Cl2 N9 O26 S
- CI COM

PAGE 1-A

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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ANSWER 10 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
L8
    256350-18-2 REGISTRY
RN
    Entered STN: 17 Feb 2000
ED
    Vancomycin, N3''-[(4'-chloro[1,1'-biphenyl]-4-yl)methyl]-6'-deoxy-6'-
CN
     [[(2,4,6-trimethylphenyl)sulfonyl]oxy]-, mono(trifluoroacetate) (salt)
     (9CI) (CA INDEX NAME)
    STEREOSEARCH
FS
    C88 H94 Cl3 N9 O26 S . C2 H F3 O2
MF
SR
LC
     STN Files:
                 CA, CAPLUS, CASREACT, USPATFULL
     СМ
          1
     CRN 256350-17-1
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Absolute stereochemistry.

C88 H94 Cl3 N9 O26 S

ОН

PAGE 3-B

CM 2

CRN 76-05-1 CMF C2 H F3 O2

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 11 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256350-17-1 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, N3''-[(4'-chloro[1,1'-biphenyl]-4-yl)methyl]-6'-deoxy-6'[[(2,4,6-trimethylphenyl)sulfonyl]oxy]- (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C88 H94 Cl3 N9 O26 S
- CI COM
- SR CA
- LC STN Files: CA, CAPLUS, USPAT2, USPATFULL

Absolute stereochemistry.

PAGE 1-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1907 TO DATE)
2 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L8 ANSWER 12 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN

RN 256350-16-0 REGISTRY

ED Entered STN: 17 Feb 2000

CN Vancomycin, 6'-deoxy-6'-[[(2,4,6-trimethylphenyl)sulfonyl]oxy]- (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C75 H85 Cl2 N9 O26 S

CI COM

SR CA

LC STN Files: CA, CAPLUS, CASREACT, USPAT7ULL

Absolute stereochemistry.

ОН

_NH2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 2 REFERENCES IN FILE CA (1907 TO DATE)
- 2 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- L8 ANSWER 13 OF 13 REGISTRY COPYRIGHT 2007 ACS on STN
- RN 256349-88-9 REGISTRY
- ED Entered STN: 17 Feb 2000
- CN Vancomycin, N3'',56-bis[(2-propenyloxy)carbonyl]-, 2-propenyl ester, 6'-(2,4,6-trimethylbenzenesulfonate) (9CI) (CA INDEX NAME)
- FS STEREOSEARCH
- MF C86 H97 Cl2 N9 O30 S
- SR CA
- LC STN Files: CA, CAPLUS, CASREACT, USPAT2, USPATFULL

Absolute stereochemistry.

PAGE 1-B

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 3 REFERENCES IN FILE CA (1907 TO DATE)
- 3 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> d his

(FILE 'HOME' ENTERED AT 15:55:45 ON 12 APR 2007)

FILE 'REGISTRY' ENTERED AT 15:55:53 ON 12 APR 2007

STRUCTURE UPLOADED L1

0 S L1 SSS SAM L2

0 S L1 SSS FUL L3

L4STRUCTURE UPLOADED

L5 12 S L4 SSS SAM

STRUCTURE UPLOADED L6

L70 S L6 SSS SAM

13 S L6 SSS FUL L8

=> file uspatfull caplus COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

374.95 375.16

FULL ESTIMATED COST

FILE 'USPATFULL' ENTERED AT 16:07:16 ON 12 APR 2007 CA INDEXING COPYRIGHT (C) 2007 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'CAPLUS' ENTERED AT 16:07:16 ON 12 APR 2007 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2007 AMERICAN CHEMICAL SOCIETY (ACS)

=> s 113

L13 NOT FOUND

The L-number entered has not been defined in this session, or it has been deleted. To see the L-numbers currently defined in this session, enter DISPLAY HISTORY at an arrow prompt (=>).

=> s 18

8 L8 L9

=> dup rem 19

PROCESSING COMPLETED FOR L9

8 DUP REM L9 (0 DUPLICATES REMOVED) L10

=> d l10 bib ab 1-8

ANSWER 1 OF 8 USPATFULL on STN T.10

2005:88018 USPATFULL AN

Glycopeptide antibiotics, combinatorial libraries of glycopeptide ΤI antibiotics and methods of producing same

Kahne, Daniel, Princeton, NJ, UNITED STATES IN

Kerns, Robert, Troy, MI, UNITED STATES

Fukuzawa, Seketsu, Tokyo, JAPAN

Ge, Min, Princeton, NJ, UNITED STATES

Thompson, Christopher, Milford, MA, UNITED STATES

Trustees of Princeton University (U.S. corporation) PA

рT

US 2005075483 A1 20050407 US 2003-676391 A1 20031001 (10) AΤ

Division of Ser. No. US 1999-353368, filed on 14 Jul 1999, GRANTED, Pat. RLI No. US 6710168

PRAT US 1999-134839P 19990519 (60)

DT Utility FS APPLICATION

LREP WOODCOCK WASHBURN LLP, ONE LIBERTY PLACE, 46TH FLOOR, 1650 MARKET STREET, PHILADELPHIA, PA, 19103

CLMN Number of Claims: 116

ECL Exemplary Claim: 1

DRWN 26 Drawing Page(s)

LN.CNT 4349

CAS INDEXING IS AVAILABLE FOR THIS PATENT.

AB A glycopeptide of the formula A.sub.1-A.sub.2-A.sub.3-A.sub.4-A.sub.5-A.sub.6-A.sub.7, in which each dash represents a covalent bond; wherein A.sub.1 comprises a modified or unmodified α-amino acid residue, alkyl, aryl, aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclic, heterocyclic-carbonyl, heterocyclic-alkyl, heterocyclic-alkyl-carbonyl, alkylsulfonyl, arylsulfonyl, guanidinyl, carbamoyl, or xanthyl; wherein each of A.sub.2 to A.sub.7 comprises a modified or unmodified α-amino acid residue, whereby (i) A.sub.1 is linked to an amino group on A.sub.2, (ii) each of A.sub.2, A.sub.4 and A.sub.6 bears an aromatic side chain, which aromatic side chains are cross-linked together by two or more covalent bonds, and (iii) A.sub.7 bears a terminal carboxyl, ester, amide, or N-substituted amide group;

and wherein one or more of A.sub.1 to A.sub.7 is linked via a glycosidic bond to one or more glycosidic groups each having one or more sugar residues, at least one of the sugar residues bearing one or more substituents of the formula YXR, N.sup.+(R.sub.1).dbd.CR.sub.2R.sub.3, N.dbd.PR.sub.1R.sub.2R.sub.3, N.sup.+R.sub.1R.sub.2R.sub.3 or P.sup.+R.sub.1R.sub.2R.sub.3 in which Y is a single bond, O, NR.sub.1 or S; X is O, NR.sub.1, S, SO.sub.2, C(O)O, C(O)S, C(S)O, C(S)S, C(NR.sub.1)O, C(O)NR.sub.1, or halo (in which case Y and R are absent).

A chemical library comprising a plurality of the glycopeptides of the invention.

A method for preparing a glycopeptide by glycosylation of an aglycone derived from a glycopeptide antibiotic.

A method for preparing a glycopeptide by preparing a pseudoaglycone from a glycopeptide antibiotic and glycosylating the pseudoaglycone.

L10 ANSWER 2 OF 8 USPATFULL on STN

AN 2004:139596 USPATFULL

TI Glycopeptide antibiotics, combinatorial libraries of glycopeptide antibiotics and methods of producing same

IN Kahne, Daniel, Princeton, NJ, UNITED STATES

Kerns, Robert, Troy, MI, UNITED STATES

Fukuzawa, Seketsu, Tokyo, JAPAN

Ge, Min, Princeton, NJ, UNITED STATES

Thompson, Christopher, Milford, MA, UNITED STATES

PI US 2004106772 A1 20040603

AI US 2003-631883 Al 20030731 (10)

RLI Division of Ser. No. US 1999-353368, filed on 14 Jul 1999, GRANTED, Pat. No. US 6710168

PRAI US 1999-134839P 19990519 (60)

DT Utility

FS APPLICATION

LREP WOODCOCK WASHBURN LLP, ONE LIBERTY PLACE, 46TH FLOOR, 1650 MARKET STREET, PHILADELPHIA, PA, 19103

CLMN Number of Claims: 116

ECL Exemplary Claim: 1

DRWN 26 Drawing Page(s)

LN.CNT 4343

CAS INDEXING IS AVAILABLE FOR THIS PATENT.

AB A glycopeptide of the formula A.sub.1-A.sub.2-A.sub.3-A.sub.4-A.sub.5-A.sub.6-A.sub.7, in which each dash represents a covalent bond; wherein A.sub.1 comprises a modified or unmodified α -amino acid residue,

alkyl, aryl, aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclic, heterocyclic-carbonyl, heterocyclic-alkyl, heterocyclic-alkyl-carbonyl, alkylsulfonyl, arylsulfonyl, guanidinyl, carbamoyl, or xanthyl; wherein each of A.sub.2 to A.sub.7 comprises a modified or unmodified α -amino acid residue, whereby (i) A.sub.1 is linked to an amino group on A.sub.2, (ii) each of A.sub.2, A.sub.4 and A.sub.6 bears an aromatic side chain, which aromatic side chains are cross-linked together by two or more covalent bonds, and (iii) A.sub.7 bears a terminal carboxyl, ester, amide, or N-substituted amide group;

and wherein one or more of A.sub.1 to A.sub.7 is linked via a glycosidic bond to one or more glycosidic groups each having one or more sugar residues, at least one of the sugar residues bearing one or more substituents of the formula YXR, N.sup.+(R.sub.1)=CR.sub.2R.sub.3, N=PR.sub.1R.sub.2R.sub.3, N.sup.+R.sub.1R.sub.2R.sub.3 or P.sup.+R.sub.1R.sub.2R.sub.3 in which Y is a single bond, O, NR, or S; X is O, NR.sub.1, S, SO.sub.2, C(O)O, C(O)S, C(S)O, C(S)S, C(NR.sub.1)O, C(O)NR.sub.1, or halo (in which case Y and R are absent).

A chemical library comprising a plurality of the glycopeptides of the invention.

A method for preparing a glycopeptide by glycosylation of an aglycone derived from a glycopeptide antibiotic.

A method for preparing a glycopeptide by preparing a pseudoaglycone from a glycopeptide antibiotic and glycosylating the pseudoaglycone.

```
L10 ANSWER 3 OF 8 USPATFULL on STN
        2004:72659 USPATFULL
AN
        Glycopeptide antibiotics, combinatorial libraries of glycopeptide
TI
        antibiotics and methods of producing same
        Kahne, Daniel, Princeton, NJ, United States
ΙN
        Kerns, Robert, Troy, MI, United States
        Fukuzawa, Seketsu, Tokyo, JAPAN
        Ge, Min, Prineton, NJ, United States
        Thompson, Christopher, Milford, MA, United States
        The Trustees of the University of Princeton, Princeton, NJ, United
 PA
        States (U.S. corporation)
 PΤ
        US 6710168
                           B1 20040323
                                19990714 (9)
        US 1999-353368
ΑI
                            19990519 (60)
        US 1999-134839P
 PRAI
        Utility
DT
        GRANTED
 FS
 EXNAM Primary Examiner: Celsa, Bennett
        Woodcock Washburn LLP
 LREP
        Number of Claims: 20
 CLMN
        Exemplary Claim: 1
 ECL
        26 Drawing Figure(s); 26 Drawing Page(s)
 DRWN
 LN.CNT 4017
· CAS INDEXING IS AVAILABLE FOR THIS PATENT.
        A glycopeptide of the formula A.sub.1--A.sub.2--A.sub.3--A.sub.4--
 AB
        bond; wherein A.sub.1 comprises a modified or unmodified \alpha-amino
```

A glycopeptide of the formula A.sub.1--A.sub.2--A.sub.3--A.sub.4--A.sub.5--A.sub.6--A.sub.7, in which each dash represents a covalent bond; wherein A.sub.1 comprises a modified or unmodified α-amino acid residue, alkyl, aryl, aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclic, heterocyclic-carbonyl, heterocyclic-alkyl, heterocyclic-alkyl-carbonyl, alkylsulfonyl, arylsulfonyl, guanidinyl, carbamoyl, or xanthyl; wherein each of A.sub.2 to A.sub.7 comprises a modified or unmodified α-amino acid residue, whereby (i) A.sub.1 is linked to an amino group on A.sub.2, (ii) each of A.sub.2, A.sub.4 and A.sub.6 bears an aromatic side chain, which aromatic side chains are cross-linked together by two or more covalent bonds, and (iii) A.sub.7 bears a terminal carboxyl, ester, amide, or N-substituted amide group;

and wherein one or more of A.sub.1 to A.sub.7 is linked via a glycosidic

bond to one or more glycosidic groups each having one or more sugar residues, at least one of the sugar residues bearing one or more substituents of the formula YXR, N.sup.+(R.sub.1).dbd.CR.sub.2R.sub.3, N.dbd.PR.sub.1R.sub.2R.sub.3, N.sup.+R.sub.1R.sub.2R.sub.3 or P.sup.+R.sub.1R.sub.2R.sub.3 in which Y is a single bond, O, NR.sub.1 or S; X is O, NR.sub.1, S, SO.sub.2, C(O)O, C(O)S, C(S)O, C(S)S, C(NR.sub.1)O, C(O)NR.sub.1, or halo (in which case Y and R are absent).

A chemical library comprising a plurality of the glycopeptides of the invention.

A method for preparing a glycopeptide by glycosylation of an aglycone derived from a glycopeptide antibiotic.

A method for preparing a glycopeptide by preparing a pseudoaglycone from a glycopeptide antibiotic and glycosylating the pseudoaglycone.

```
L10 ANSWER 4 OF 8 USPATFULL on STN
ΑN
       2003:319226 USPATFULL
       Glycopeptide antibacterial compounds, compositions containing same and
TI
       methods of using same
       Kim, Ronald M., Hoboken, NJ, UNITED STATES
IN
       Kahne, Daniel E., Princeton, NJ, UNITED STATES
       Chapman, Kevin T., Scotch Plains, NJ, UNITED STATES
       Princeton University, Princeton, NJ, UNITED STATES (U.S. corporation)
PA
                           A1 20031204
PΙ
       US 2003224975
                           B2 20050111
       US 6841661
       US 2002-262858
                           A1 20021003 (10)
ΑI
       Continuation of Ser. No. US 2000-574225, filed on 19 May 2000, GRANTED,
RLI
       Pat. No. US 6498238
       US 1999-134841P
                           19990519 (60)
PRAI
DT
       Utility
FS
       APPLICATION
       Robert L. Price, McDermott, Will & Emery, 600 13th Street, N.W.,
LREP
       Washington, DC, 20005-3096
       Number of Claims: 47
CLMN
       Exemplary Claim: 1
ECL
       No Drawings
DRWN
LN.CNT 1877
CAS INDEXING IS AVAILABLE FOR THIS PATENT.
       The present invention relates to vancomycin analogs in which the
AΒ
       vancosamine residue is substituted with a lipid-like substituent that
       includes a first aryl moiety and a second aryl moiety joined together by
       a flexible linker moiety, that is not a single bond directly joining the
       first aryl moiety and the second aryl moiety, and a glucose C-6
       substituent modified to be other than the naturally occurring hydroxyl
       group, or pharmaceutically acceptable salts thereof.
L10 ANSWER 5 OF 8 USPATFULL on STN
       2002:340371 USPATFULL
AN
       Glycopeptide antibacterial compounds, compositions containing same and
TT
       methods of using same
       Kim, Ronald M., Hoboken, NJ, United States
TN
       Kahne, Daniel E., Princeton, NJ, United States
       Chapman, Kevin T., Scotch Plains, NJ, United States
       Princeton University, Princeton, NJ, United States (U.S. corporation)
PΑ
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       US 6498238
       US 2000-574225
ΑI
                               20000519 (9)
       US 1999-134841P
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      Primary Examiner: Riley, Jezia
EXNAM
LREP
       McDermott, Will & Emery
CLMN
       Number of Claims: 47
ECL
       Exemplary Claim: 1
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0 Drawing Figure(s); 0 Drawing Page(s) DRWN LN.CNT 1575 CAS INDEXING IS AVAILABLE FOR THIS PATENT. The present invention relates to vancomycin analogs in which the AB vancosamine residue is substituted with a lipid-like substituent that includes a first aryl moiety and a second aryl moiety joined together by a flexible linker moiety, that is not a single bond directly joining the first aryl moiety and the second aryl moiety, and a glucose C-6 substituent modified to be other than the naturally occurring hydroxyl group, or pharmaceutically acceptable salts thereof. ANSWER 6 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN 2001:798248 CAPLUS AN DN 135:331681 Preparation of glycopeptide vancomycin analogs as antibacterial agents ΤI IN Kahne, Daniel; Walker, Suzanne Trustees of Princeton University, USA PΑ PCT Int. Appl., 68 pp. SO CODEN: PIXXD2 DT Patent LA English FAN.CNT 2 APPLICATION NO. DATE PATENT NO. KIND DATE ______ ____ _____ WO 2001-US11040 20010405 A2 20011101 PΙ WO 2001081372 A3 20020516 WO 2001081372 W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG 20020411 US 2001-818787 20010328 US 2002042365 A1 US 6699836 B2 20040302 A5 .20011107 AU 2001-51321 20010405 AU 2001051321 P PRAI US 2000-199382P 20000425 A US 2001-818787 20010328 P 19990402 US 1999-127516P W 20010405 WO 2001-US11040 MARPAT 135:331681 OS Antibacterial glycopeptide vancomycin analogs I bearing optional AB modifications to the C6 position of the glucose residue attached to the amino acid four of the vancomycin heptapeptide chain, wherein X is O, S, substituted amine; R is H, substituted alkyl, aryl, aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclic, heterocyclic-carbonyl, heterocyclic-alkyl-carbonyl, alkylsulfonyl, arylsulfonyl, aminoalkyl; R1 and R2 are independently substituted alkyl, aryl, aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclic, heterocyclic-carbonyl, heterocyclic-alkyl-carbonyl, alkylsulfonyl, arylsulfonyl; substituted alkyl, acyl, CHO; Z is N3, functionalized O, N, S atoms; are disclosed. Methods of making the compds. and methods of using the compds. to treat a bacterial infection in a host are also disclosed. Thus, I [X = NH, Z = OH, R = CH2CH2CH2CH2NH2, R1 = H, R2 = 4-[(3,4dichlorophenyl) methoxy] benzyl] was prepared and tested in vitro for their antibacterial activity (MIC = $0.25-6.25 \mu g/mL$). ANSWER 7 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN L10 2000:824286 CAPLUS AN DN 134:5162 TI Preparation of glycopeptides as antibacterial agents

Kim, Ronald M.; Kahne, Daniel E.; Chapman, Kevin T.

Merck & Co., Inc., USA; Princeton University

IN

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PCT Int. Appl., 89 pp.
SO
     CODEN: PIXXD2
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PRAI US 1999-134841P
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     US 2000-574225
OS
     MARPAT 134:5162
     Glycopeptides I [R is a polar substituent; K-Arl-Z-Ar2 is a lipid-like
AΒ
     substituent where Arl and Ar2 are aromatic or heterocyclic groups, each
     optionally substituted with R1 [R1 = halo, R2, CN, NO2, CF3,
     fluoromethoxy, NHSO2R2, OR2, SR2, NR22, N+R23, C(O)NR22, SO2NR22,
     heterocyclyl, CO2R2, C(O)R2, OC(O)R2, NR2C(O)R2, or NHC(O)R2; R2 = H,
     aryl, alkyl, arylalkyl, (heterocyclyl)alkyl, aroyl, alkanoyl, alkanoyloxy,
     alkanoylamino, alkylsulfonyl, arylsulfonyl; two R2 groups may form one or
     more aromatic or heterocyclic rings]; K and Z are carbonyl, sulfonyl,
     alkylene, alkyleneoxy, oxyalkylene, alkyleneamino, aminoalkylene,
     alkyleneoxyalkylene, alkylenethio, thioalkylene, alkylenecarbonyl,
     aminocarbonyl or carbonylamino, alkyleneaminocarbonyl,
     aminocarbonylalkylene, O, O2C, CO2, alkylene, alkyleneoxycarbonyl,
     oxycarbonylalkylene, aminosulfonyl or sulfonylamino; Z is not a singe
     bond] were prepared as antibacterial agents. Thus, N-[4-(3,4-
     dichlorobenzyloxy)benzyl]-N-glucose-C6-amino-vancomycin, prepared from
     vancomycin hydrochloride by a multistep sequence involving condensation
     with 4-(3,4-dichlorobenzyloxy) benzaldehyde, showed MIC = 0.125 \mu g/mL
     against Staphylococcus aureus Septicemia (in vivo).
               THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
RE.CNT 3
               ALL CITATIONS AVAILABLE IN THE RE FORMAT
     ANSWER 8 OF 8 CAPLUS COPYRIGHT 2007 ACS on STN
L10
     2000:68479 CAPLUS
AN
DN
     132:122934
     Preparation of glycopeptide antibiotics and their combinatorial libraries
ΤI
     Kahne, Daniel; Kerns, Robert; Fukuzawa, Seketsu; Ge, Min; Thompson,
IN
     Christopher
     Princeton University, USA
PΑ
SO
     PCT Int. Appl., 159 pp.
     CODEN: PIXXD2
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     Patent
     English
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PRAI US 1998-150690P
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os
     Glycopeptides A1-A2-A3-A4-A5-A6-A7 [A1 comprises a modified or unmodified
AB
     α-amino acid residue, alkyl, aryl, aralkyl, alkanoyl, aroyl,
     aralkanoyl, heterocyclyl, heterocyclylcarbonyl, heterocyclylalkyl,
     heterocyclylalkylcarbonyl, alkylsulfonyl, arylsulfonyl, guanidinyl,
     carbamoyl, or xanthyl; each of A2 to A7 comprises a modified or unmodified
     \alpha\text{-amino} acid residue, where (i) Al is linked to an amino group on
     A2, (ii) each of A2, A4 and A6 bears an aromatic side chain which is
     cross-linked by two or more covalent bonds, and (iii) A7 bears a terminal
     carboxyl, ester, amide, or N-substituted amide group; one or more of Al to
     A7 is linked via a glycosidic bond to one or more glycosidic groups each
     having one or more sugar residues, at least one of the sugar residues
     bearing one or more substituents of the formula YXR, N+R1:CR2R3,
     N:PR1R2R3, N+R1R2R3 or P+R1R2R3 in which Y is a single bond, O, NR1 or S;
     X is O, NR1, S, SO2, C(O)O, C(O)S, C(S)O, C(S)S, C(NR1)O, C(O)NR1, or halo
     (in which case Y and R are absent); R, R1, R2, and R3 are H, alkyl, aryl,
     aralkyl, alkanoyl, aroyl, aralkanoyl, heterocyclyl, heterocyclylcarbonyl,
     heterocyclylalkyl, heterocyclylalkylcarbonyl, alkylsulfonyl, or
     arylsulfonyl] and their pharmaceutically acceptable salts or a chemical
     library comprising a plurality of the glycopeptides of the invention were
     prepared for use as antibiotics. Thus, glucose-C6 modified vancomycin
     derivs. were prepared and assayed for antimicrobial activity (min.
     inhibitory concns. are tabulated).
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THERE ARE 8 CITED REFERENCES AVAILABLE FOR THIS RECORD

ALL CITATIONS AVAILABLE IN THE RE FORMAT

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RE.CNT 8

SINCE FILE TOTAL ENTRY SESSION 24.57 399.73

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

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SESSION WILL BE HELD FOR 120 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 16:11:18 ON 12 APR 2007